C:\Program Files\Stnexp\Queries\10673521-3.str

$$0^{e_{-H}}$$
 $0^{e_{-H}}$ $0^{$

chain nodes:

7 8 9 10 11 12 13 14 15 16 17 18 19 20 21 23 30 31 32 33 34 35 36 37 38 57 58

ring nodes:

1 2 3 4 5 6 39 40 41 42 43 44 45 46 47 48 49

chain bonds:

2-7 4-30 6-57 7-58 8-9 10-11 11-12 13-14 14-15 14-16 17-18 19-20 19-21 21-23 31-32 32-33 34-35 36-37 37-38

ring bonds:

1-2 1-6 2-3 3-4 4-5 5-6 39-40 39-43 40-41 41-42 42-43 44-45 44-49 45-46 46-47 47-48 48-49

exact/norm bonds:

2-7 4-30 6-57 7-58 8-9 10-11 11-12 13-14 14-15 14-16 17-18 19-20 19-21 21-23 31-32 32-33 34-35 36-37 37-38 39-40 39-43 40-41 41-42 42-43 44-45 44-49 45-46 46-47 47-48 48-49

normalized bonds:

1-2 1-6 2-3 3-4 4-5 5-6

G1:H,Cy,Ak

G2:[*1],[*2],[*3],[*4],[*5]

G3:0,S,N

G4:[*6],[*7],[*8],[*9],[*10]

Match level:

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:CLASS8:CLASS9:CLASS10:CLASS11:CLASS 12:CLASS13:CLASS14:CLASS15:CLASS16:CLASS17:CLASS18:CLASS19:CLASS20:CLASS21:CLASS 23:CLASS30:CLASS31:CLASS32:CLASS33:Atom 34:CLASS35:Atom 36:CLASS37:CLASS38:CLASS 39:Atom 40:Atom 41:Atom 42:Atom 43:Atom 44:Atom 45:Atom 46:Atom 47:Atom 48:Atom 49:Atom 57:CLASS8:CLASS

Generic attributes:

58:

Saturation

: Unsaturated